



Full wwPDB EM Validation Report ⓘ

Jun 30, 2026 – 12:48 PM JST

PDB ID : 9WXH / pdb_00009wxh
EMDB ID : EMD-66346
Title : Cryo-EM structure of the type III-D2 CRISPR-Cas effector complex bound to a cognate target RNA in the pre-cleavage state
Authors : Mitsuda, Y.; Ishikawa, J.; Nagahata, N.; Hiraizumi, M.; Yamashita, K.; Nishimasu, H.
Deposited on : 2025-09-25
Resolution : 2.60 Å(reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

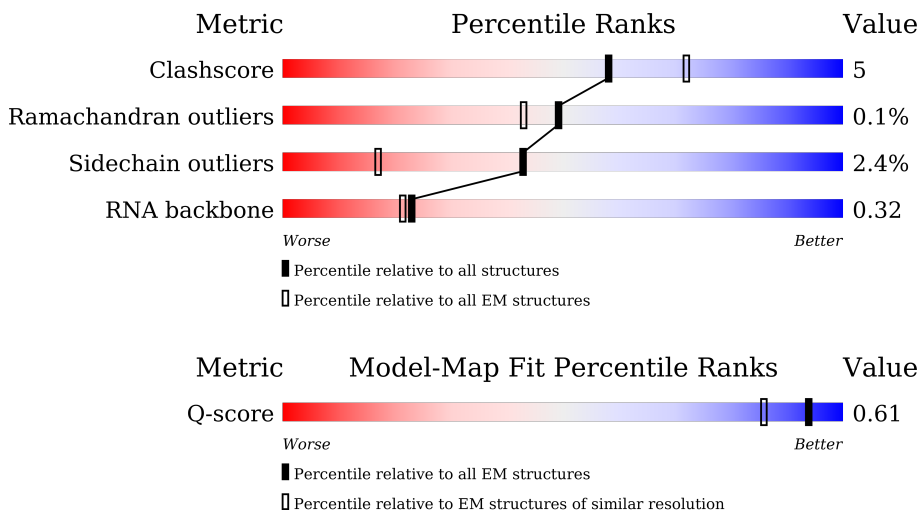
EMDB validation analysis : 0.0.1.dev133
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 2.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



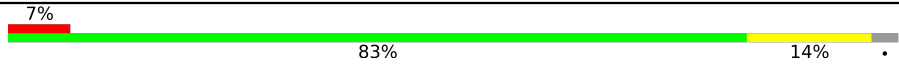

Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	8728 (2.10 - 3.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	43	
2	B	592	
3	C	635	

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Mol	Chain	Length	Quality of chain
4	D	1235	
5	E	47	

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 20118 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a RNA chain called CRISPR RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
1	A	40	841	378	141	283	39	0	0

- Molecule 2 is a protein called Cas10.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	494	4012	2532	699	769	12	0	0

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	MET	-	initiating methionine	UNP A0A9D0P2E0
B	-18	GLY	-	expression tag	UNP A0A9D0P2E0
B	-17	SER	-	expression tag	UNP A0A9D0P2E0
B	-16	SER	-	expression tag	UNP A0A9D0P2E0
B	-15	HIS	-	expression tag	UNP A0A9D0P2E0
B	-14	HIS	-	expression tag	UNP A0A9D0P2E0
B	-13	HIS	-	expression tag	UNP A0A9D0P2E0
B	-12	HIS	-	expression tag	UNP A0A9D0P2E0
B	-11	HIS	-	expression tag	UNP A0A9D0P2E0
B	-10	HIS	-	expression tag	UNP A0A9D0P2E0
B	-9	SER	-	expression tag	UNP A0A9D0P2E0
B	-8	SER	-	expression tag	UNP A0A9D0P2E0
B	-7	GLY	-	expression tag	UNP A0A9D0P2E0
B	-6	LEU	-	expression tag	UNP A0A9D0P2E0
B	-5	VAL	-	expression tag	UNP A0A9D0P2E0
B	-4	PRO	-	expression tag	UNP A0A9D0P2E0
B	-3	ARG	-	expression tag	UNP A0A9D0P2E0
B	-2	GLY	-	expression tag	UNP A0A9D0P2E0
B	-1	SER	-	expression tag	UNP A0A9D0P2E0
B	0	HIS	-	expression tag	UNP A0A9D0P2E0

- Molecule 3 is a protein called CRISPR type III-associated protein domain-containing protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	C	626	4912	3132	828	935	17	0	0

- Molecule 4 is a protein called TIGR03986 family CRISPR-associated RAMP protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	D	1194	9433	5911	1699	1791	32	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	0	SER	-	expression tag	UNP A0A9D0P2B2

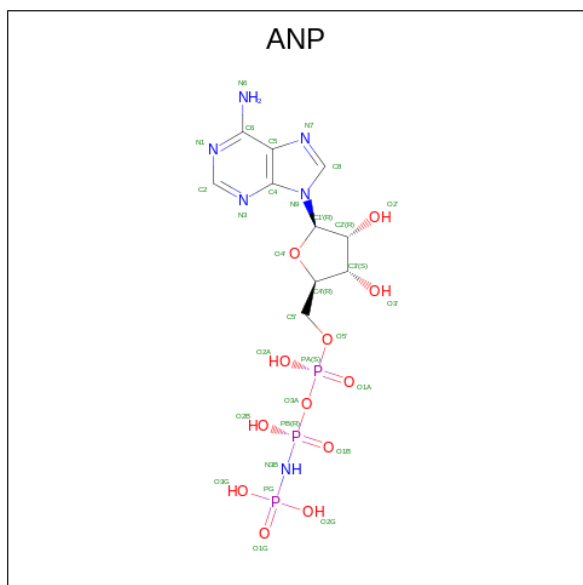
- Molecule 5 is a RNA chain called Cognate target RNA (CTR).

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
5	E	38	813	365	153	257	38	0	0

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

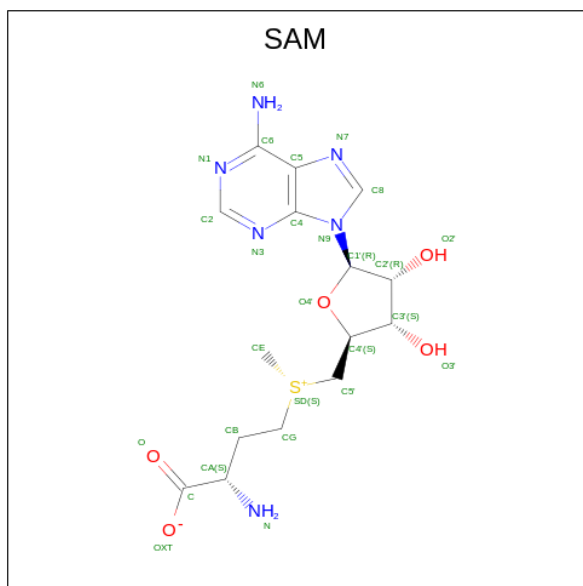
Mol	Chain	Residues	Atoms		AltConf
6	B	3	Total	Mg	0
			3	3	
6	C	1	Total	Mg	0
			1	1	

- Molecule 7 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
7	B	1	31	10	6	12	3	0

- Molecule 8 is S-ADENOSYLMETHIONINE (CCD ID: SAM) (formula: $C_{15}H_{22}N_6O_5S$) (labeled as "Ligand of Interest" by depositor).



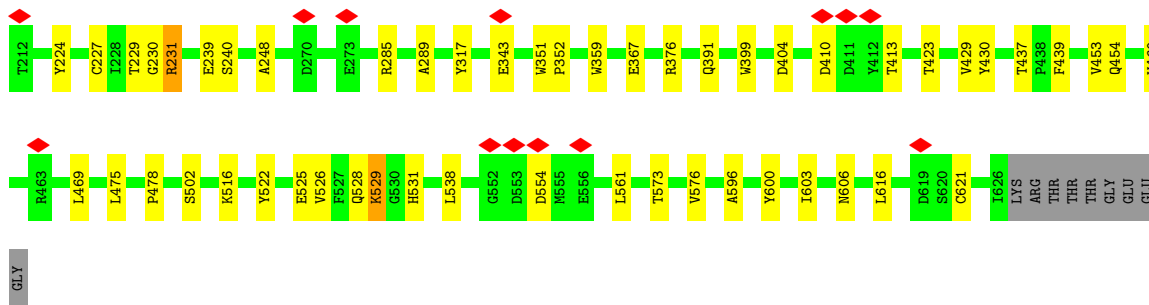
Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	S	
8	B	1	54	30	12	10	2	1

- Molecule 9 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

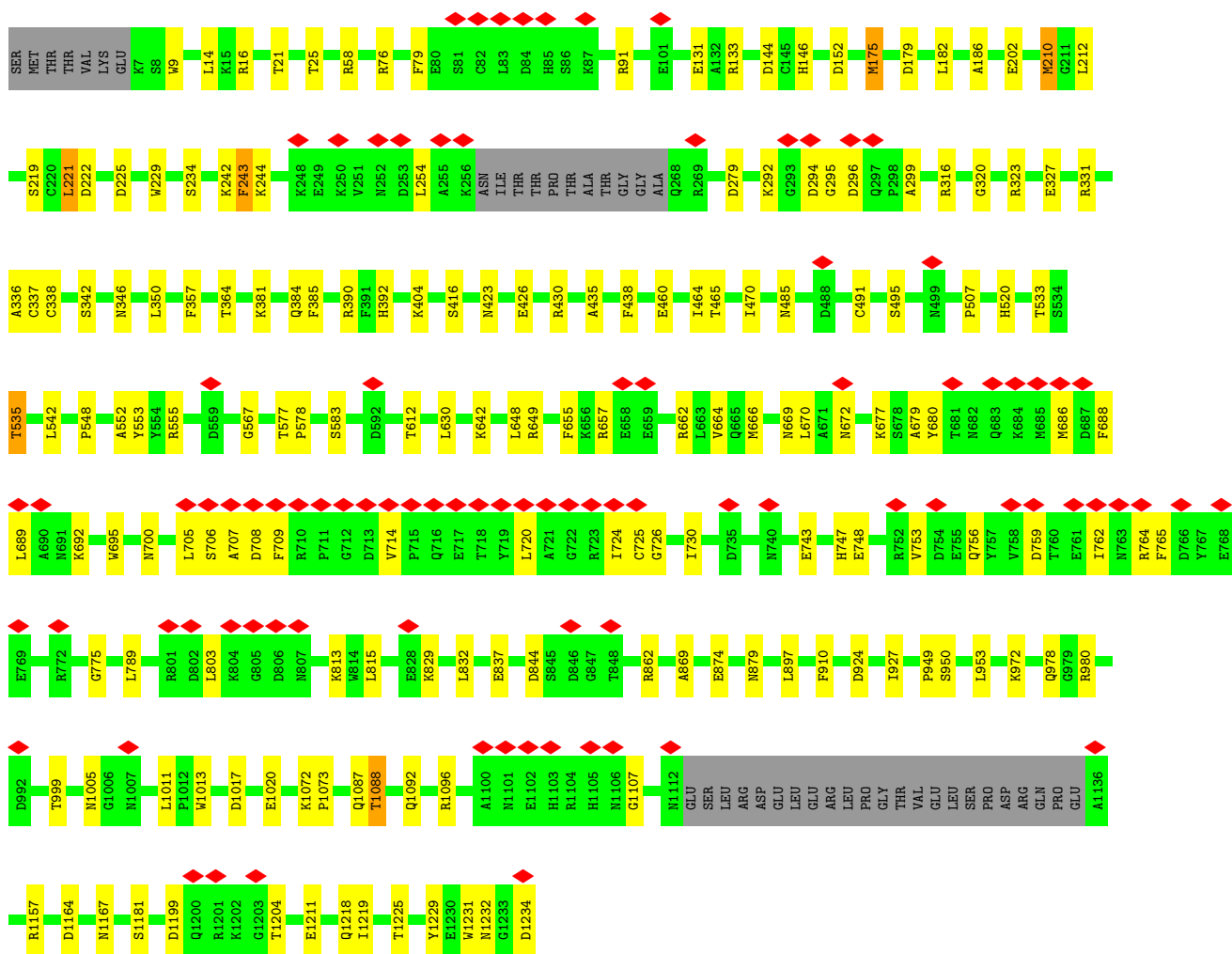
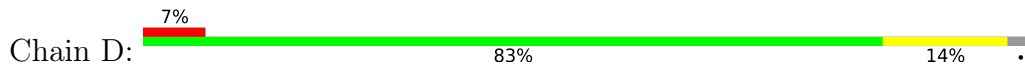
Mol	Chain	Residues	Atoms		AltConf
9	D	1	Total 1	Zn 1	0

- Molecule 10 is water.

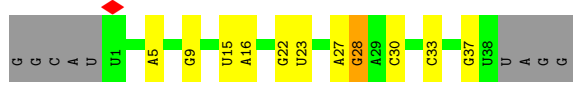
Mol	Chain	Residues	Atoms		AltConf
10	A	5	Total 5	O 5	0
10	C	6	Total 6	O 6	0
10	D	3	Total 3	O 3	0
10	E	3	Total 3	O 3	0



• Molecule 4: TIGR03986 family CRISPR-associated RAMP protein



• Molecule 5: Cognate target RNA (CTR)



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	371927	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	49.833	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	1.531	Depositor
Minimum map value	-0.849	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.030	Depositor
Recommended contour level	0.111	Depositor
Map size (Å)	298.8, 298.8, 298.8	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.99599993, 0.99599993, 0.99599993	Depositor

5 Model quality i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: SAM, MG, ANP, ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.67	0/938	1.37	11/1458 (0.8%)
2	B	0.57	0/4088	0.99	5/5504 (0.1%)
3	C	0.59	0/5032	0.96	4/6845 (0.1%)
4	D	0.60	0/9648	0.98	7/13041 (0.1%)
5	E	0.63	0/911	1.12	5/1417 (0.4%)
All	All	0.60	0/20617	1.01	32/28265 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
2	B	0	1
All	All	0	2

There are no bond length outliers.

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	3	U	O3'-P-O5'	-10.40	88.39	104.00
4	D	210	MET	CG-SD-CE	-8.05	83.18	100.90
2	B	307	ASP	CA-CB-CG	7.55	120.15	112.60
4	D	243	PHE	N-CA-CB	7.15	120.62	109.69
4	D	465	THR	CA-CB-OG1	-6.81	99.39	109.60
4	D	553	TYR	CB-CA-C	-6.60	97.97	109.86
5	E	27	A	O3'-P-O5'	-6.49	94.26	104.00
3	C	97	TYR	N-CA-CB	6.41	120.78	110.65
1	A	31	U	C4'-C3'-O3'	6.30	118.85	109.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	23	G	C4'-C3'-O3'	-6.14	103.79	113.00
3	C	430	TYR	N-CA-CB	-6.09	100.29	111.37
1	A	8	C	C3'-C2'-O2'	-6.01	105.58	114.60
3	C	97	TYR	CB-CA-C	-5.91	101.23	110.74
5	E	33	C	O3'-P-O5'	-5.79	95.31	104.00
5	E	23	U	O3'-P-O5'	-5.77	95.35	104.00
1	A	35	G	O3'-P-O5'	-5.62	95.57	104.00
1	A	23	G	O3'-P-O5'	-5.62	95.58	104.00
4	D	152	ASP	CA-CB-CG	5.52	118.12	112.60
1	A	26	A	C3'-C2'-O2'	-5.49	106.36	114.60
1	A	18	A	C1'-C2'-O2'	5.41	116.51	108.40
3	C	475	LEU	N-CA-CB	-5.37	102.16	110.65
1	A	27	A	C1'-O4'-C4'	-5.31	104.39	109.70
4	D	980	ARG	CD-NE-CZ	5.19	131.66	124.40
1	A	9	U	C1'-O4'-C4'	-5.16	104.55	109.70
5	E	9	G	O3'-P-O5'	-5.12	96.33	104.00
2	B	558	TYR	CA-C-N	-5.11	116.32	122.35
2	B	558	TYR	C-N-CA	-5.11	116.32	122.35
2	B	449	ARG	NE-CZ-NH2	5.09	123.78	119.20
2	B	405	HIS	CA-CB-CG	-5.07	108.73	113.80
1	A	5	A	O3'-P-O5'	-5.05	96.43	104.00
4	D	175	MET	CG-SD-CE	-5.04	89.81	100.90
5	E	30	C	O3'-P-O5'	-5.01	96.48	104.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	39	G	Sidechain
2	B	244	VAL	Peptide

5.2 Too-close contacts [\(i\)](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	841	0	427	6	0
2	B	4012	0	3971	42	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	4912	0	4864	50	0
4	D	9433	0	9215	93	0
5	E	813	0	412	2	0
6	B	3	0	0	0	0
6	C	1	0	0	0	0
7	B	31	0	13	0	0
8	B	54	0	44	1	0
9	D	1	0	0	0	0
10	A	5	0	0	0	0
10	C	6	0	0	0	0
10	D	3	0	0	0	0
10	E	3	0	0	0	0
All	All	20118	0	18946	184	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (184) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:435:ALA:HB2	4:D:470:ILE:HG21	1.47	0.96
4:D:79:PHE:HE2	4:D:186:ALA:HB2	1.36	0.86
4:D:79:PHE:CE2	4:D:186:ALA:HB2	2.17	0.78
4:D:1092:GLN:OE1	4:D:1096:ARG:NH1	2.16	0.78
2:B:165:PRO:HB3	2:B:174:ARG:HD3	1.67	0.76
3:C:317:TYR:HB3	3:C:616:LEU:HD21	1.67	0.76
4:D:649:ARG:O	4:D:670:LEU:HG	1.87	0.75
2:B:39:ILE:HD11	2:B:153:TYR:CE2	2.24	0.73
2:B:123:SER:OG	8:B:605[B]:SAM:N7	2.21	0.73
4:D:495:SER:O	4:D:535:THR:OG1	2.07	0.73
4:D:662:ARG:NH2	4:D:775:GLY:O	2.23	0.71
2:B:248:PHE:CE1	2:B:252:ASN:ND2	2.62	0.67
4:D:655:PHE:HA	4:D:837:GLU:HG3	1.78	0.66
4:D:677:LYS:HA	4:D:747:HIS:ND1	2.10	0.66
2:B:20:ASN:O	2:B:405:HIS:CE1	2.50	0.65
1:A:3:U:O2'	3:C:231:ARG:HB2	1.96	0.65
4:D:726:GLY:HA2	4:D:753:VAL:HG22	1.79	0.65
3:C:12:LEU:CD2	3:C:188:VAL:HG22	2.27	0.64
4:D:705:LEU:O	4:D:724:ILE:N	2.26	0.64
2:B:393:SER:OG	2:B:430:ILE:HG12	1.99	0.62
4:D:144:ASP:OD1	4:D:144:ASP:C	2.42	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:343:LEU:HD12	2:B:344:TRP:N	2.15	0.61
4:D:294:ASP:O	4:D:296:ASP:N	2.33	0.61
4:D:1204:THR:O	4:D:1232:ASN:ND2	2.33	0.61
3:C:53:GLN:HG3	4:D:9:TRP:CD2	2.36	0.60
3:C:204:PRO:HA	3:C:423:THR:HG21	1.81	0.60
2:B:205:ASN:C	2:B:205:ASN:OD1	2.45	0.60
4:D:507:PRO:O	4:D:1157:ARG:NH2	2.33	0.60
4:D:730:ILE:HA	4:D:748:GLU:OE1	2.02	0.59
4:D:803:LEU:HD21	4:D:813:LYS:HB2	1.85	0.59
1:A:27:A:O2'	1:A:28:U:OP2	2.22	0.57
3:C:367:GLU:OE2	3:C:391:GLN:NE2	2.38	0.57
2:B:236:ARG:O	2:B:240:ARG:HG2	2.04	0.57
4:D:202:GLU:OE2	4:D:316:ARG:NH2	2.37	0.57
4:D:655:PHE:CE2	4:D:664:VAL:HG23	2.41	0.56
3:C:137:ILE:HD13	3:C:180:MET:HE1	1.87	0.56
4:D:688:PHE:CG	4:D:714:VAL:HG13	2.41	0.56
3:C:162:THR:O	3:C:166:HIS:ND1	2.34	0.56
3:C:317:TYR:HB3	3:C:616:LEU:CD2	2.35	0.55
4:D:323:ARG:HG2	4:D:327:GLU:OE1	2.06	0.55
2:B:324:GLN:OE1	2:B:360:TRP:HZ2	1.88	0.55
3:C:289:ALA:HB1	3:C:399:TRP:HB3	1.87	0.55
4:D:426:GLU:HB3	4:D:430:ARG:NH1	2.21	0.55
4:D:385:PHE:CE1	4:D:612:THR:HG21	2.42	0.55
4:D:680:TYR:HA	4:D:686:MET:HG2	1.89	0.55
4:D:700:ASN:ND2	4:D:764:ARG:HG3	2.21	0.55
4:D:336:ALA:N	4:D:485:ASN:O	2.38	0.55
3:C:73:TRP:CE3	3:C:165:LYS:HD3	2.42	0.54
3:C:169:THR:HG23	4:D:234:SER:OG	2.07	0.54
4:D:58:ARG:NH2	4:D:133:ARG:O	2.41	0.54
2:B:324:GLN:OE1	2:B:360:TRP:CZ2	2.61	0.54
4:D:346:ASN:HA	4:D:350:LEU:HD21	1.89	0.53
2:B:444:ASN:ND2	2:B:447:ASP:OD2	2.41	0.53
3:C:46:LEU:C	3:C:46:LEU:CD1	2.82	0.53
4:D:655:PHE:HE2	4:D:664:VAL:HG23	1.74	0.53
4:D:789:LEU:HD21	4:D:862:ARG:HB3	1.89	0.53
4:D:657:ARG:HB3	4:D:662:ARG:HD2	1.91	0.52
4:D:520:HIS:CE1	4:D:1225:THR:HG22	2.44	0.52
4:D:1011:LEU:HD13	4:D:1013:TRP:CZ2	2.45	0.52
4:D:357:PHE:O	4:D:364:THR:HB	2.10	0.52
2:B:394:SER:HA	2:B:431:ALA:O	2.10	0.51
5:E:15:U:O2	5:E:15:U:H2'	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:135:VAL:HG22	5:E:28:G:C6	2.44	0.51
4:D:533:THR:HG21	4:D:874:GLU:HG2	1.93	0.51
4:D:79:PHE:HE2	4:D:186:ALA:CB	2.15	0.51
4:D:1218:GLN:HG3	4:D:1219:ILE:HD12	1.93	0.51
4:D:219:SER:HB3	4:D:243:PHE:HB3	1.93	0.51
3:C:596:ALA:HA	3:C:600:TYR:O	2.11	0.50
3:C:616:LEU:N	3:C:616:LEU:HD22	2.26	0.50
4:D:91:ARG:NH1	4:D:179:ASP:OD2	2.44	0.50
4:D:1072:LYS:N	4:D:1073:PRO:HD2	2.27	0.49
2:B:155:TYR:OH	2:B:230:ASP:OD2	2.26	0.49
3:C:522:TYR:CE1	3:C:526:VAL:HG21	2.47	0.49
4:D:222:ASP:OD1	4:D:225:ASP:N	2.44	0.49
2:B:294:GLN:HB3	2:B:306:PHE:CE2	2.47	0.49
3:C:138:GLU:HG2	3:C:140:PRO:HD3	1.95	0.49
4:D:648:LEU:HD22	4:D:844:ASP:HB2	1.93	0.49
2:B:316:GLU:OE1	2:B:372:GLN:NE2	2.46	0.48
3:C:6:LEU:HD22	3:C:164:LEU:HD22	1.94	0.48
4:D:229:TRP:HB2	4:D:242:LYS:HD3	1.96	0.48
4:D:390:ARG:HB2	4:D:630:LEU:HD23	1.96	0.48
2:B:118:ILE:HD11	2:B:128:GLU:HB2	1.95	0.48
3:C:478:PRO:HB3	3:C:538:LEU:HD22	1.95	0.47
4:D:385:PHE:HE1	4:D:612:THR:HG21	1.78	0.47
4:D:706:SER:O	4:D:709:PHE:HD2	1.98	0.47
3:C:53:GLN:HA	4:D:9:TRP:CD1	2.49	0.47
2:B:460:LEU:N	2:B:461:PRO:CD	2.78	0.47
2:B:259:LYS:O	2:B:336:ARG:HA	2.15	0.47
4:D:144:ASP:OD1	4:D:146:HIS:N	2.41	0.47
4:D:1231:TRP:NE1	4:D:1234:ASP:OD2	2.45	0.47
4:D:707:ALA:O	4:D:708:ASP:HB2	2.15	0.47
2:B:364:ARG:NH1	2:B:368:LEU:HD21	2.29	0.46
3:C:163:ILE:O	3:C:167:ILE:HG12	2.16	0.46
4:D:491:CYS:SG	4:D:548:PRO:HB3	2.55	0.46
4:D:924:ASP:O	4:D:927:ILE:HG13	2.15	0.46
3:C:224:TYR:HB2	3:C:437:THR:HG23	1.98	0.46
3:C:410:ASP:OD1	3:C:413:THR:CB	2.64	0.46
2:B:169:GLU:HG2	2:B:173:SER:HB3	1.97	0.46
4:D:292:LYS:HA	4:D:299:ALA:HB2	1.98	0.46
4:D:1219:ILE:HD12	4:D:1219:ILE:N	2.31	0.45
2:B:491:TYR:CD1	2:B:553:MET:HE2	2.52	0.45
3:C:410:ASP:OD1	3:C:413:THR:OG1	2.32	0.45
4:D:279:ASP:HB3	4:D:460:GLU:HB2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:743:GLU:C	4:D:743:GLU:OE1	2.59	0.45
4:D:423:ASN:HD22	4:D:426:GLU:HG2	1.82	0.45
4:D:175:MET:HB2	4:D:179:ASP:HB2	1.97	0.45
4:D:384:GLN:NE2	4:D:583:SER:O	2.50	0.45
4:D:1219:ILE:O	4:D:1219:ILE:HG22	2.16	0.45
2:B:289:LYS:HE2	2:B:425:LYS:HG2	1.98	0.44
3:C:525:GLU:O	3:C:529:LYS:HB3	2.18	0.44
4:D:79:PHE:CE2	4:D:186:ALA:CB	2.95	0.44
3:C:46:LEU:C	3:C:46:LEU:HD12	2.43	0.44
4:D:1211:GLU:HB2	4:D:1229:TYR:OH	2.18	0.44
4:D:338:CYS:HB2	4:D:342:SER:O	2.18	0.44
2:B:479:LYS:HA	2:B:482:GLN:HG2	2.00	0.44
4:D:438:PHE:CG	4:D:464:ILE:HD13	2.52	0.44
2:B:368:LEU:HA	2:B:371:GLU:HG2	1.99	0.44
4:D:21:THR:OG1	4:D:212:LEU:HD11	2.18	0.44
4:D:381:LYS:O	4:D:404:LYS:HA	2.18	0.43
4:D:759:ASP:CG	4:D:762:ILE:HG22	2.43	0.43
2:B:485:PHE:CE1	2:B:512:ILE:HG21	2.52	0.43
4:D:327:GLU:HG2	4:D:337:CYS:SG	2.59	0.43
1:A:15:U:H5	3:C:118:ASP:OD2	2.02	0.43
2:B:20:ASN:O	2:B:405:HIS:NE2	2.51	0.43
3:C:285:ARG:HB3	3:C:404:ASP:HB3	1.99	0.43
4:D:1164:ASP:HB3	4:D:1167:ASN:ND2	2.34	0.43
4:D:700:ASN:OD1	4:D:765:PHE:HB3	2.17	0.43
3:C:317:TYR:CE1	3:C:621:CYS:SG	3.12	0.43
2:B:273:TYR:CD1	2:B:561:PRO:HB2	2.54	0.43
2:B:228:ILE:HG22	2:B:229:SER:O	2.18	0.43
2:B:249:TYR:O	2:B:253:THR:OG1	2.32	0.43
3:C:227:CYS:HA	3:C:240:SER:HB2	2.01	0.43
2:B:165:PRO:CB	2:B:174:ARG:HD3	2.45	0.42
2:B:284:TYR:HD1	2:B:350:ILE:HG12	1.84	0.42
2:B:555:LEU:O	2:B:556:TRP:C	2.63	0.42
3:C:376:ARG:HD2	4:D:9:TRP:CE2	2.55	0.42
3:C:118:ASP:HA	3:C:135:VAL:O	2.19	0.42
3:C:48:VAL:HG22	3:C:98:PHE:CD2	2.54	0.42
3:C:115:LEU:HB3	3:C:139:ASN:O	2.20	0.42
2:B:154:LYS:HE3	2:B:155:TYR:CZ	2.55	0.42
4:D:666:MET:HE2	4:D:666:MET:HB3	1.89	0.42
3:C:460:VAL:HG23	3:C:606:ASN:CG	2.45	0.42
4:D:577:THR:HB	4:D:578:PRO:HD2	2.02	0.42
3:C:429:VAL:HB	3:C:439:PHE:CZ	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:502:SER:HA	3:C:561:LEU:HD13	2.02	0.42
4:D:1017:ASP:OD2	4:D:1020:GLU:HB2	2.20	0.42
2:B:371:GLU:HG3	2:B:372:GLN:HG2	2.02	0.41
3:C:359:TRP:CE3	3:C:453:VAL:HG21	2.55	0.41
4:D:669:ASN:OD1	4:D:669:ASN:C	2.63	0.41
4:D:700:ASN:ND2	4:D:764:ARG:CG	2.82	0.41
3:C:230:GLY:HA2	3:C:239:GLU:O	2.20	0.41
4:D:1219:ILE:O	4:D:1219:ILE:CG2	2.68	0.41
1:A:2:U:O4'	3:C:248:ALA:HA	2.20	0.41
3:C:317:TYR:CB	3:C:616:LEU:HD21	2.46	0.41
3:C:343:GLU:CD	3:C:343:GLU:N	2.78	0.41
4:D:869:ALA:HB1	4:D:972:LYS:HD2	2.02	0.41
2:B:40:VAL:HG11	2:B:127:PHE:CE2	2.55	0.41
3:C:86:SER:HB2	3:C:90:LEU:O	2.20	0.41
3:C:573:THR:OG1	3:C:576:VAL:HG23	2.20	0.41
4:D:680:TYR:HA	4:D:686:MET:CG	2.50	0.41
4:D:949:PRO:O	4:D:950:SER:C	2.63	0.41
3:C:351:TRP:CG	3:C:352:PRO:HA	2.56	0.41
4:D:25:THR:HG22	4:D:210:MET:HE1	2.02	0.41
4:D:221:LEU:HD12	4:D:243:PHE:CD1	2.54	0.41
3:C:97:TYR:CD2	3:C:97:TYR:C	2.98	0.41
3:C:528:GLN:HA	3:C:531:HIS:HD2	1.85	0.41
2:B:421:ILE:HG22	2:B:424:PHE:HD2	1.85	0.41
3:C:410:ASP:OD1	3:C:413:THR:HB	2.21	0.41
4:D:679:ALA:HB1	4:D:689:LEU:CD1	2.51	0.41
4:D:999:THR:O	4:D:1181:SER:HA	2.21	0.41
2:B:271:ALA:HB2	2:B:356:TRP:CZ3	2.56	0.41
3:C:469:LEU:HD23	3:C:603:ILE:HG22	2.02	0.41
4:D:567:GLY:HA3	4:D:1087:GLN:O	2.20	0.41
4:D:542:LEU:HD21	4:D:910:PHE:HB2	2.04	0.40
4:D:1088:THR:HG22	4:D:1107:GLY:HA3	2.04	0.40
2:B:316:GLU:HB3	2:B:372:GLN:HE21	1.87	0.40
1:A:20:A:O4'	4:D:320:GLY:HA3	2.21	0.40
2:B:173:SER:O	2:B:177:LYS:HG3	2.21	0.40
2:B:6:ARG:CD	2:B:182:ASN:OD1	2.70	0.40
2:B:276:LEU:HD12	2:B:356:TRP:HB3	2.03	0.40
4:D:695:TRP:CZ2	4:D:765:PHE:HD2	2.39	0.40
1:A:6:A:O2'	1:A:7:A:H5'	2.21	0.40
4:D:229:TRP:HB2	4:D:242:LYS:CD	2.52	0.40
4:D:331:ARG:HD2	4:D:552:ALA:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	490/592 (83%)	480 (98%)	10 (2%)	0	100	100
3	C	624/635 (98%)	610 (98%)	13 (2%)	1 (0%)	43	66
4	D	1188/1235 (96%)	1163 (98%)	24 (2%)	1 (0%)	48	70
All	All	2302/2462 (94%)	2253 (98%)	47 (2%)	2 (0%)	49	70

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	D	295	GLY
3	C	229	THR

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	429/514 (84%)	417 (97%)	12 (3%)	38	66
3	C	533/540 (99%)	525 (98%)	8 (2%)	57	80
4	D	1005/1041 (96%)	977 (97%)	28 (3%)	38	66
All	All	1967/2095 (94%)	1919 (98%)	48 (2%)	43	70

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	135	THR

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Mol	Chain	Res	Type
2	B	191	SER
2	B	205	ASN
2	B	326	LYS
2	B	327	GLU
2	B	340	GLU
2	B	343	LEU
2	B	363	LYS
2	B	421	ILE
2	B	427	ARG
2	B	464	LEU
2	B	554	GLU
3	C	46	LEU
3	C	64	LEU
3	C	86	SER
3	C	231	ARG
3	C	454	GLN
3	C	516	LYS
3	C	529	LYS
3	C	554	ASP
4	D	14	LEU
4	D	16	ARG
4	D	76	ARG
4	D	131	GLU
4	D	182	LEU
4	D	221	LEU
4	D	244	LYS
4	D	254	LEU
4	D	392	HIS
4	D	416	SER
4	D	535	THR
4	D	555	ARG
4	D	642	LYS
4	D	672	ASN
4	D	692	LYS
4	D	720	LEU
4	D	725	CYS
4	D	756	GLN
4	D	815	LEU
4	D	829	LYS
4	D	832	LEU
4	D	879	ASN
4	D	897	LEU

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Mol	Chain	Res	Type
4	D	953	LEU
4	D	978	GLN
4	D	1005	ASN
4	D	1088	THR
4	D	1199	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	2	ASN
2	B	179	GLN
2	B	324	GLN
2	B	440	HIS
2	B	452	GLN
3	C	19	GLN
3	C	92	ASN
3	C	288	HIS
3	C	329	ASN
3	C	331	GLN
3	C	531	HIS
4	D	52	GLN
4	D	78	HIS
4	D	158	ASN
4	D	384	GLN
4	D	423	ASN
4	D	520	HIS
4	D	604	ASN
4	D	744	ASN
4	D	799	ASN
4	D	1005	ASN
4	D	1040	HIS
4	D	1044	ASN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	39/43 (90%)	13 (33%)	8 (20%)
5	E	37/47 (78%)	5 (13%)	0
All	All	76/90 (84%)	18 (23%)	8 (10%)

All (18) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	3	U
1	A	9	U
1	A	10	U
1	A	13	C
1	A	14	U
1	A	19	C
1	A	20	A
1	A	21	U
1	A	22	G
1	A	28	U
1	A	30	C
1	A	31	U
1	A	36	U
5	E	5	A
5	E	16	A
5	E	22	G
5	E	28	G
5	E	37	G

All (8) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	7	A
1	A	9	U
1	A	13	C
1	A	15	U
1	A	19	C
1	A	21	U
1	A	25	U
1	A	27	A

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 8 ligands modelled in this entry, 5 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	ANP	B	604	6	32,33,33	1.19	4 (12%)	44,52,52	1.35	4 (9%)
8	SAM	B	605[B]	6	27,29,29	0.43	0	36,42,42	0.63	1 (2%)
8	SAM	B	605[A]	-	27,29,29	0.41	0	36,42,42	0.54	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	ANP	B	604	6	-	5/18/38/38	0/3/3/3
8	SAM	B	605[B]	6	-	1/16/33/33	0/3/3/3
8	SAM	B	605[A]	-	-	1/16/33/33	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	604	ANP	PB-O1B	3.94	1.52	1.46
7	B	604	ANP	PG-O1G	3.54	1.51	1.46
7	B	604	ANP	PG-O2G	-2.50	1.50	1.56
7	B	604	ANP	PB-O2B	-2.14	1.51	1.56

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	604	ANP	O2B-PB-O1B	5.66	121.79	109.92
7	B	604	ANP	O1G-PG-N3B	-3.82	106.14	111.77
7	B	604	ANP	O1B-PB-N3B	2.51	115.47	111.77
7	B	604	ANP	O3A-PB-N3B	-2.45	99.81	106.59
8	B	605[B]	SAM	CG-SD-C5'	-2.02	98.26	103.40

There are no chirality outliers.

All (7) torsion outliers are listed below:

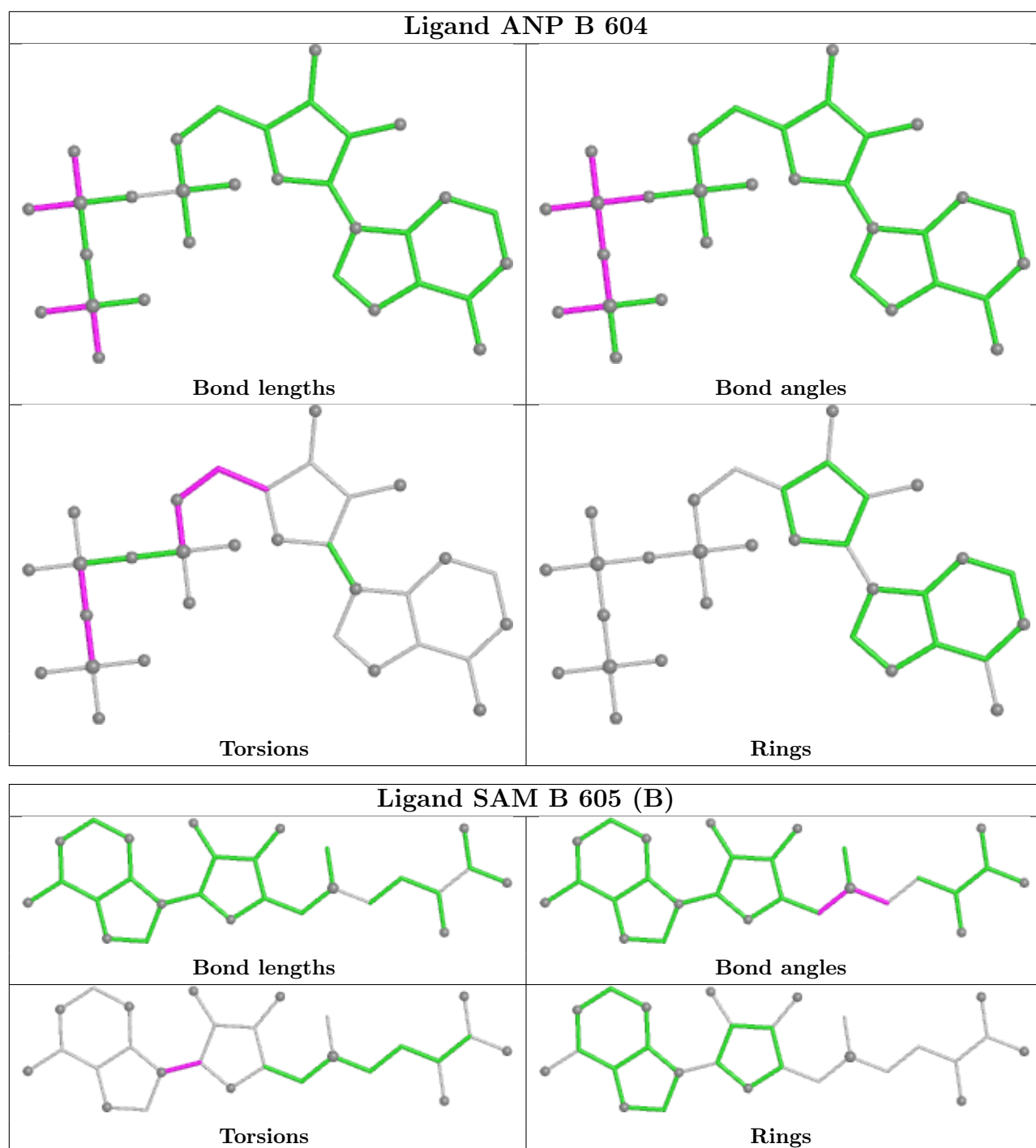
Mol	Chain	Res	Type	Atoms
7	B	604	ANP	PB-N3B-PG-O1G
7	B	604	ANP	PG-N3B-PB-O1B
8	B	605[A]	SAM	CA-CB-CG-SD
7	B	604	ANP	C4'-C5'-O5'-PA
7	B	604	ANP	O4'-C4'-C5'-O5'
8	B	605[B]	SAM	C2'-C1'-N9-C8
7	B	604	ANP	C5'-O5'-PA-O1A

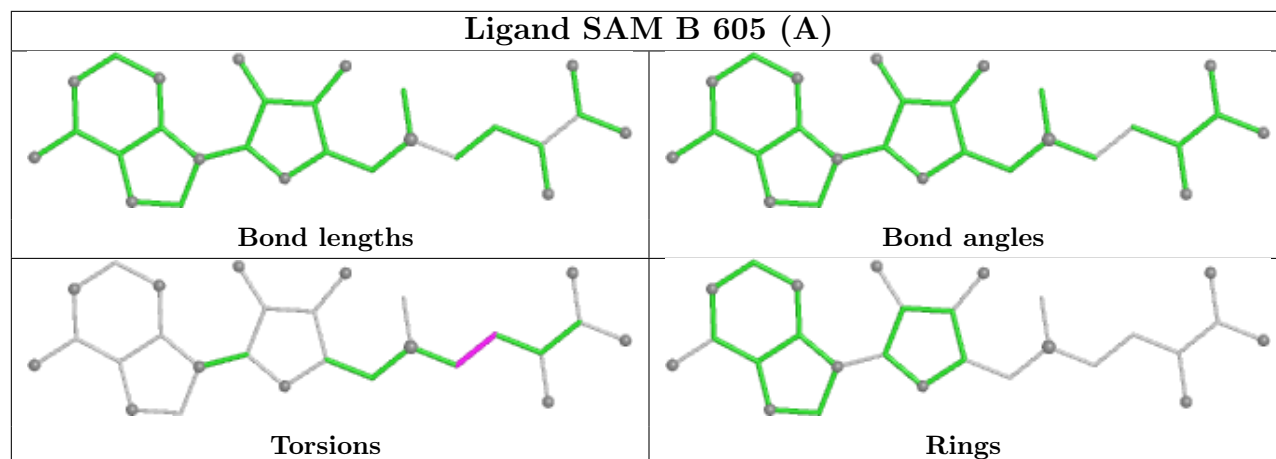
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	B	605[B]	SAM	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

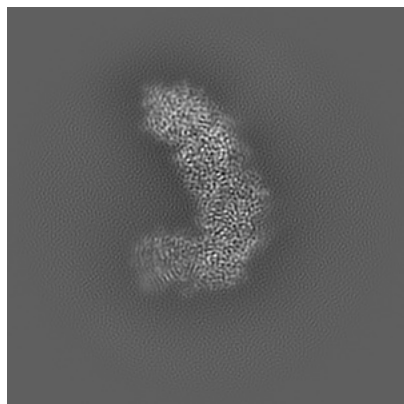
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-66346. These allow visual inspection of the internal detail of the map and identification of artifacts.

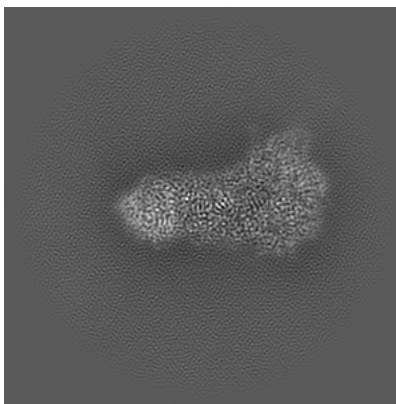
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

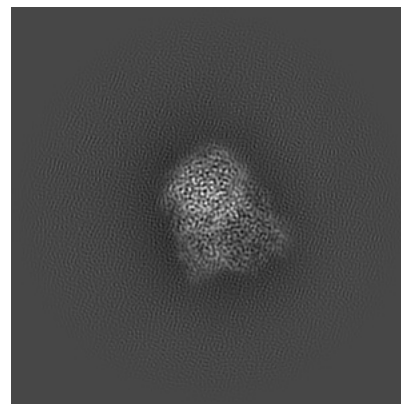
6.1.1 Primary map



X

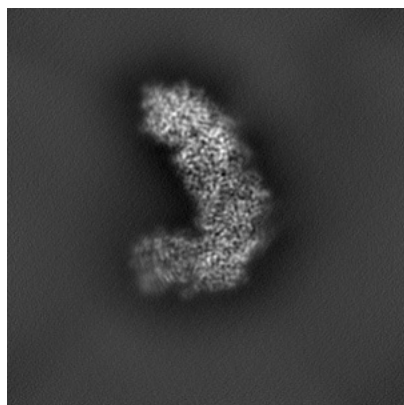


Y

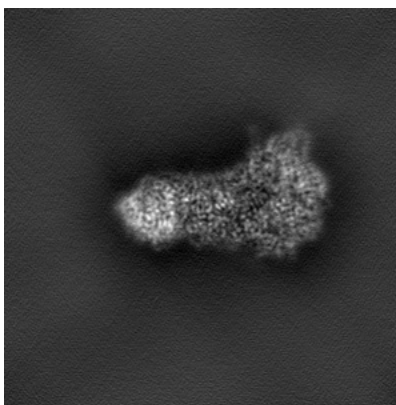


Z

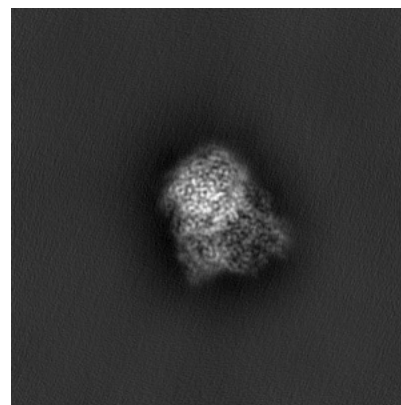
6.1.2 Raw map



X



Y

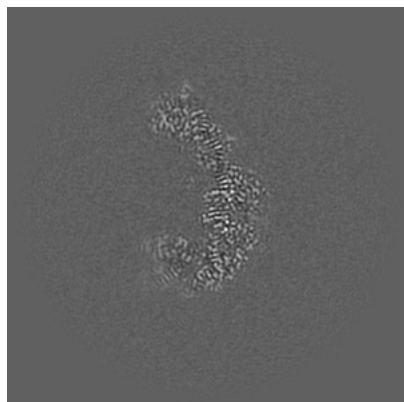


Z

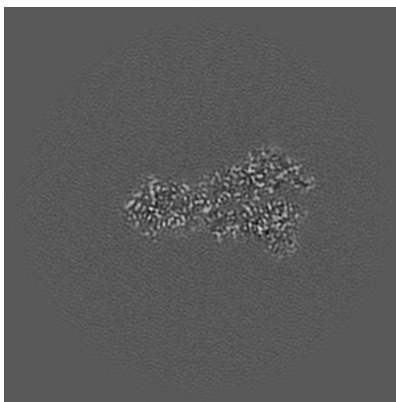
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

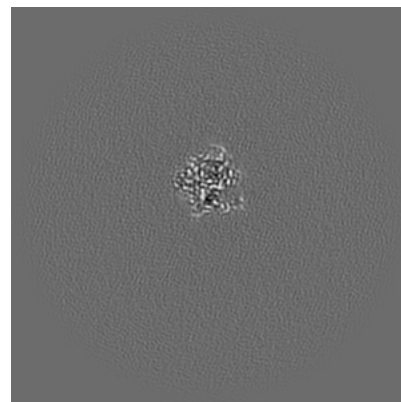
6.2.1 Primary map



X Index: 150

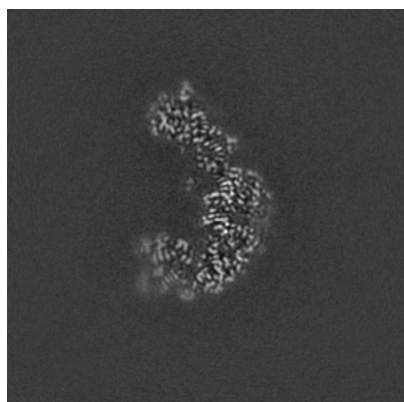


Y Index: 150

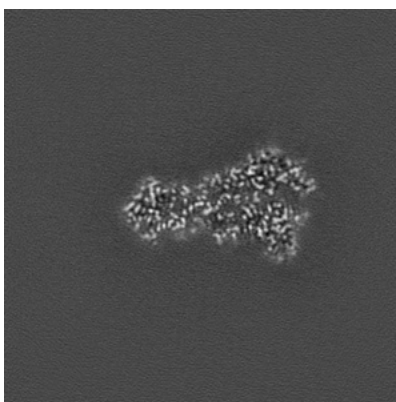


Z Index: 150

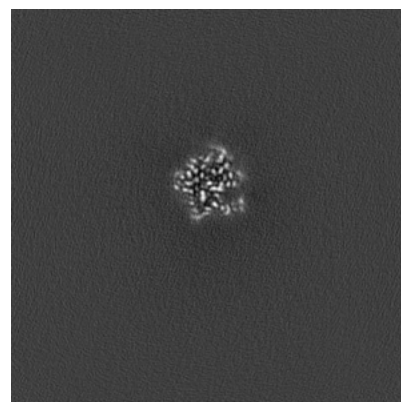
6.2.2 Raw map



X Index: 150



Y Index: 150

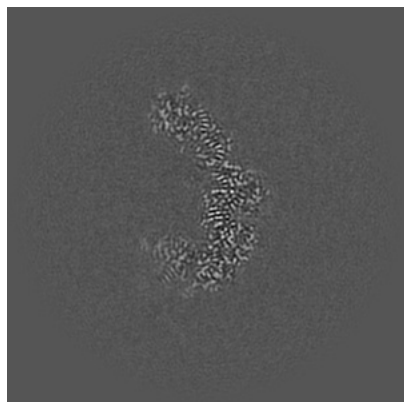


Z Index: 150

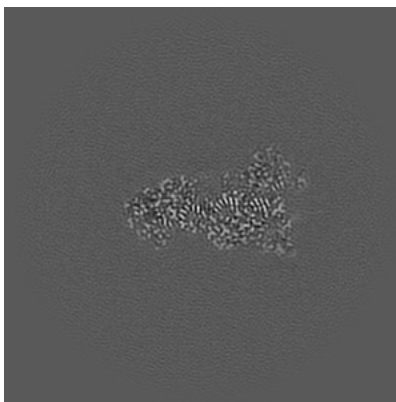
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

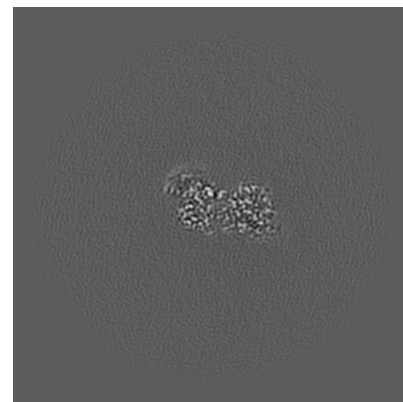
6.3.1 Primary map



X Index: 149

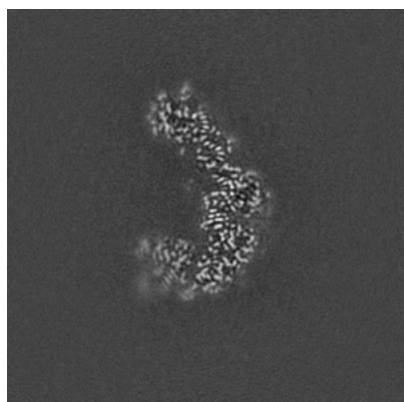


Y Index: 158

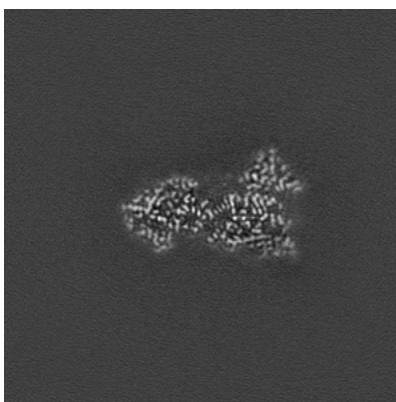


Z Index: 192

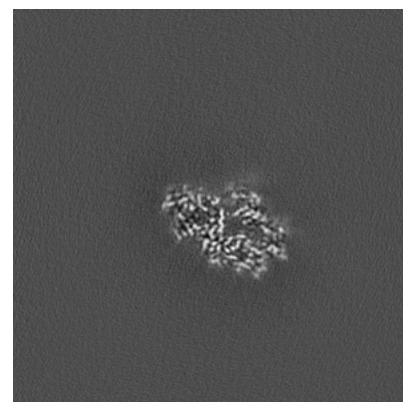
6.3.2 Raw map



X Index: 149



Y Index: 160

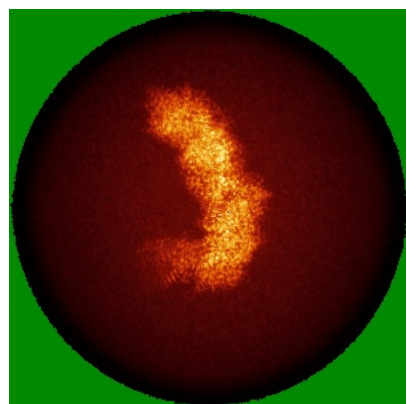


Z Index: 209

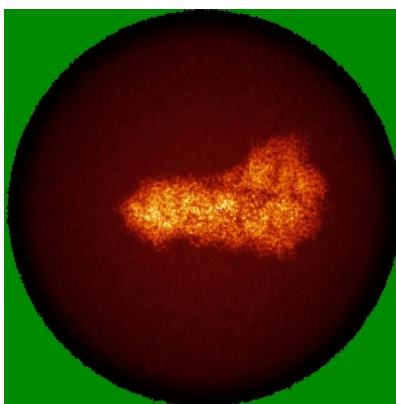
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

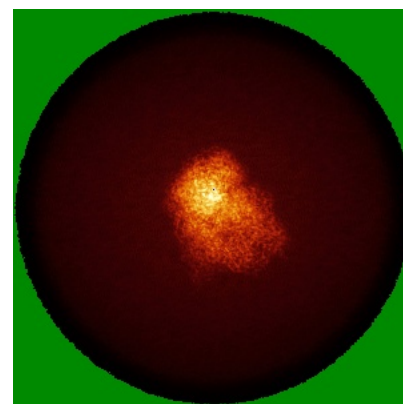
6.4.1 Primary map



X

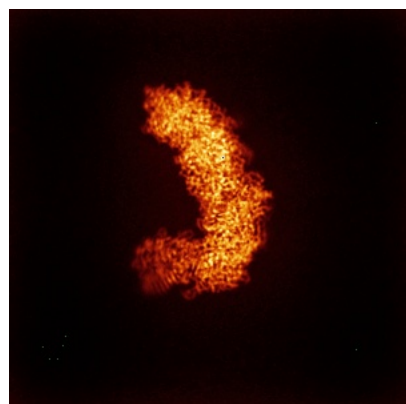


Y

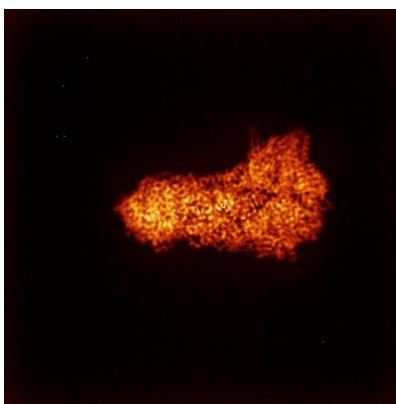


Z

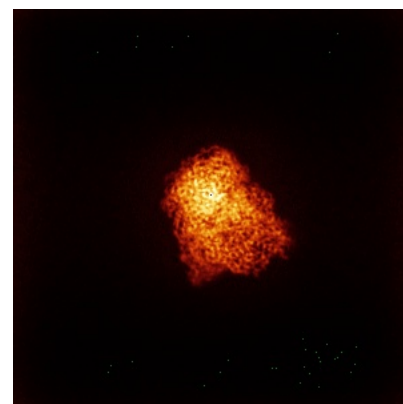
6.4.2 Raw map



X



Y

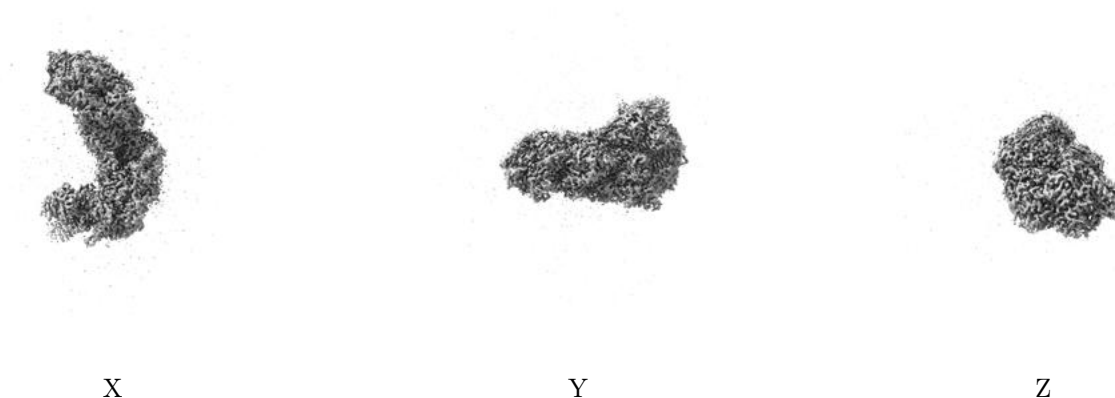


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

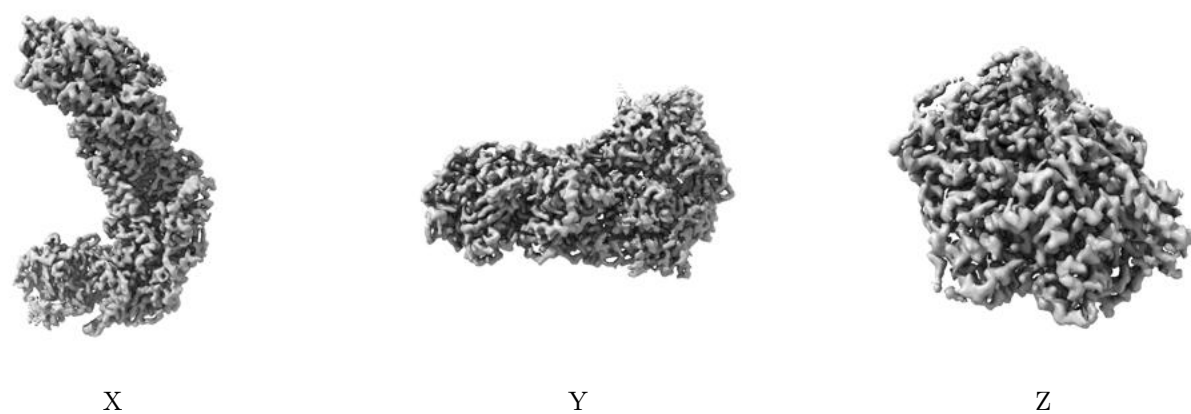
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.111. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

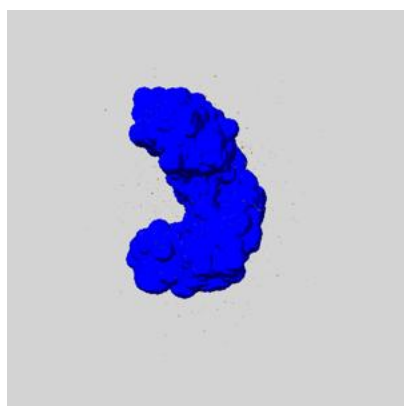
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

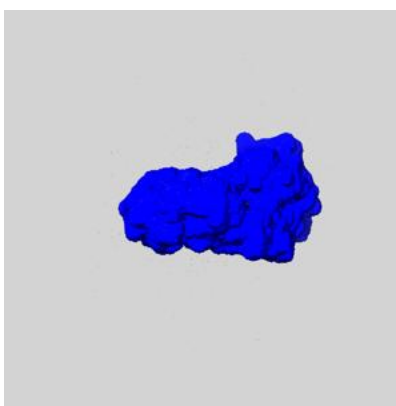
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

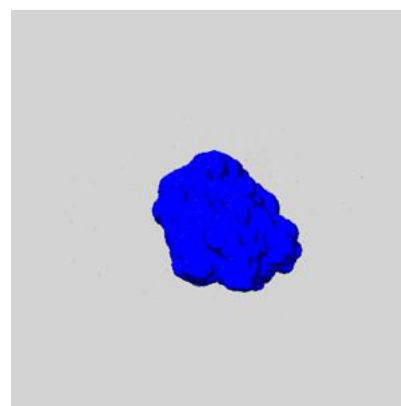
6.6.1 emd_66346_msk_1.map [i](#)



X



Y

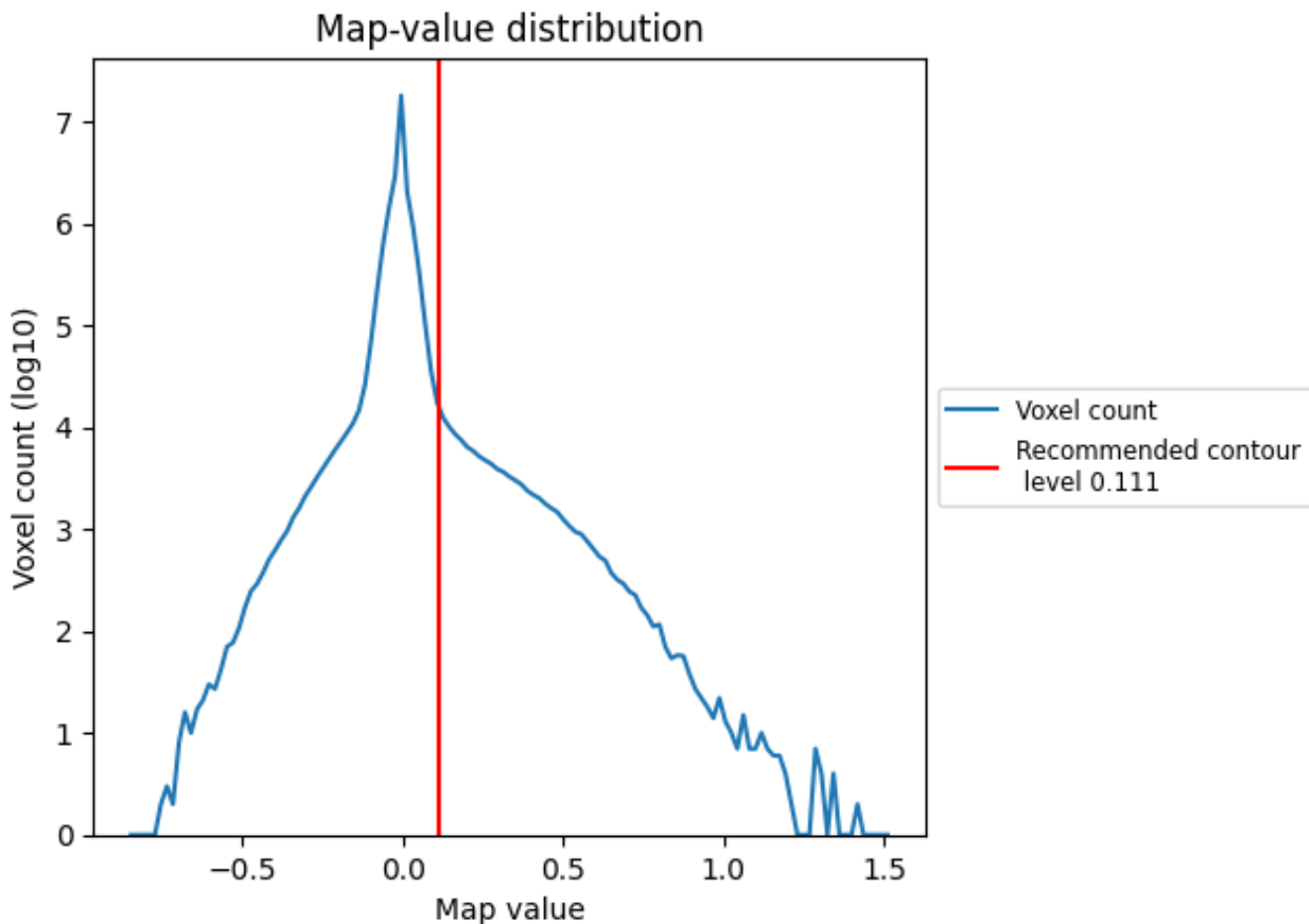


Z

7 Map analysis [i](#)

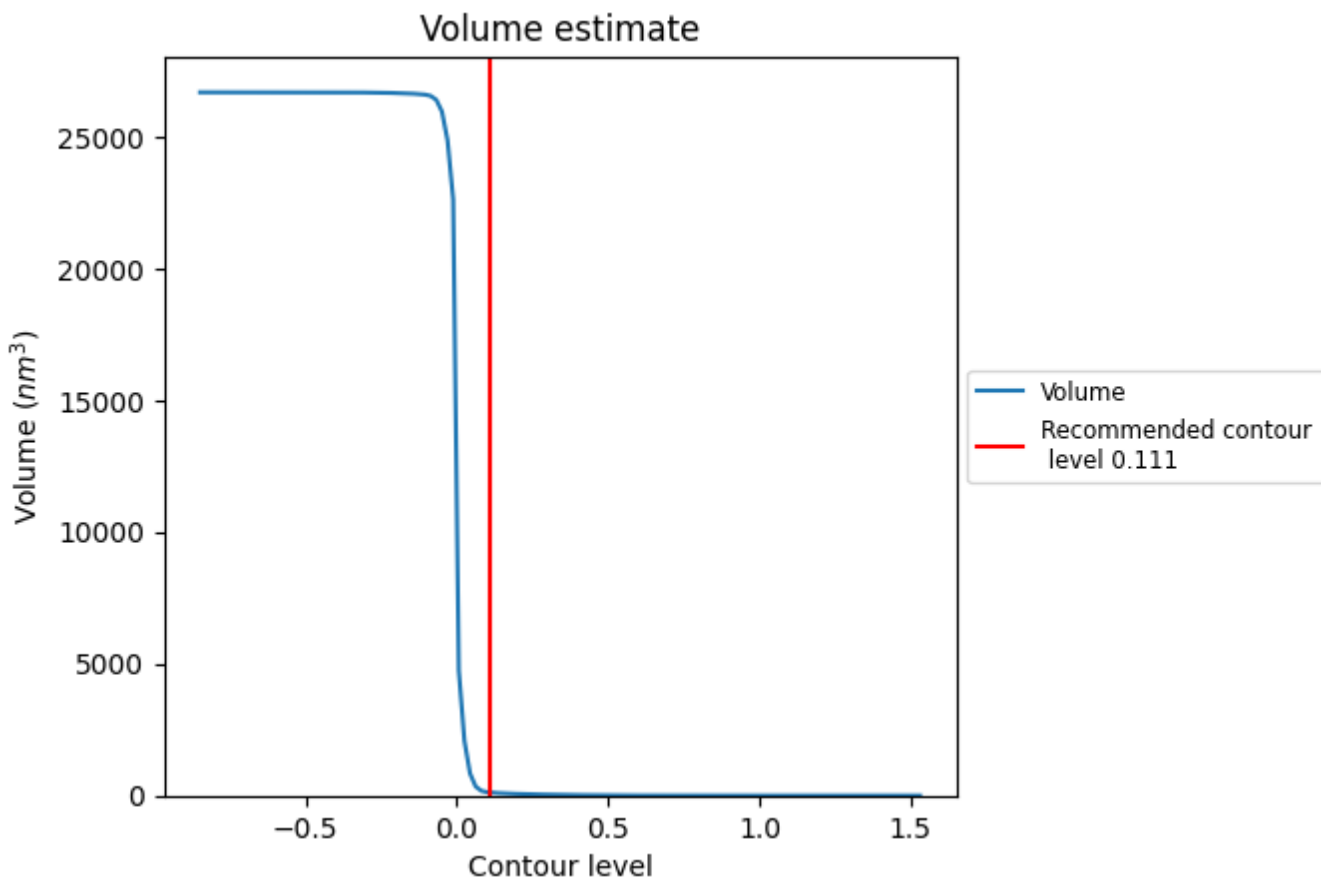
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

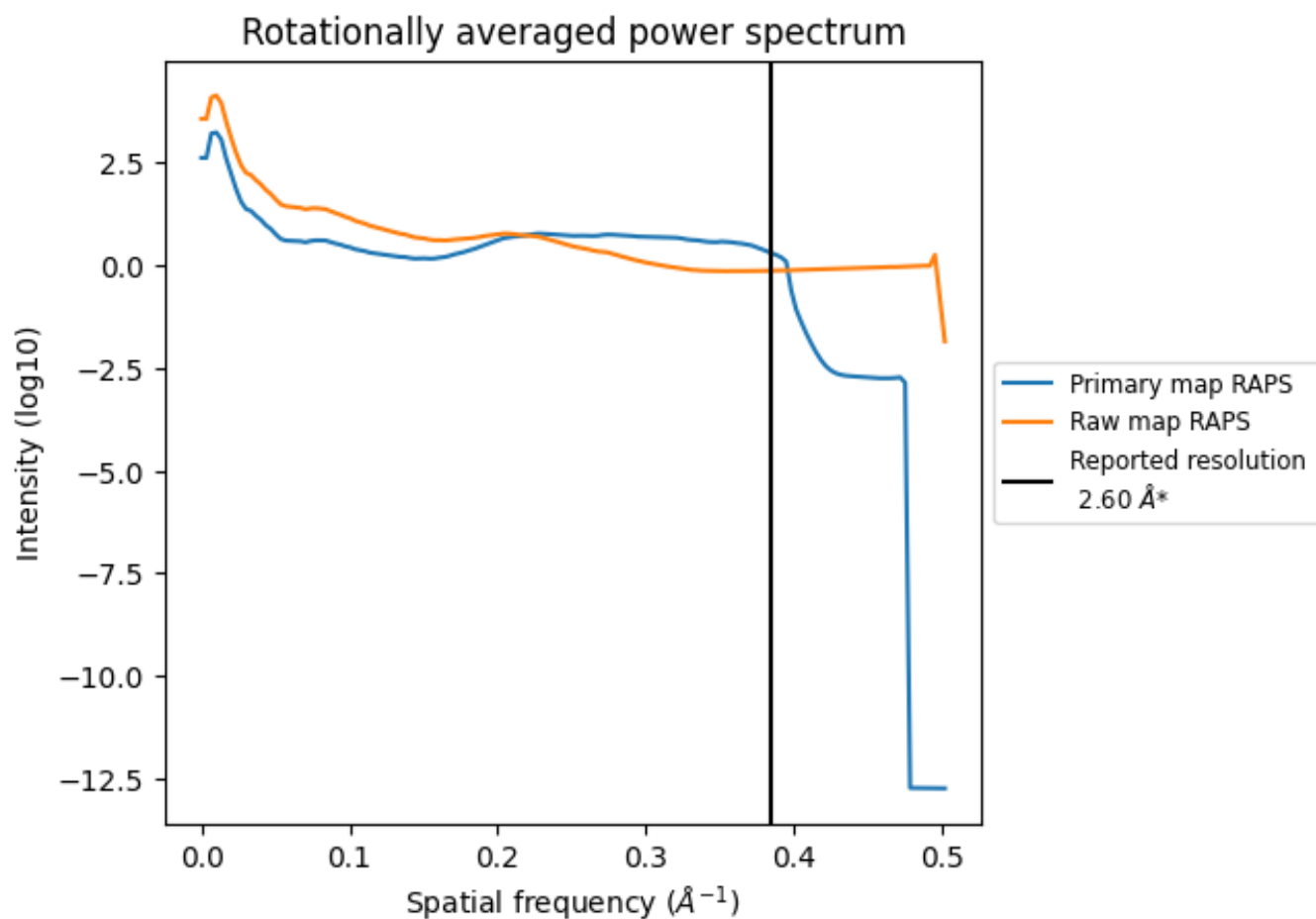
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 115 nm^3 ; this corresponds to an approximate mass of 104 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum [i](#)

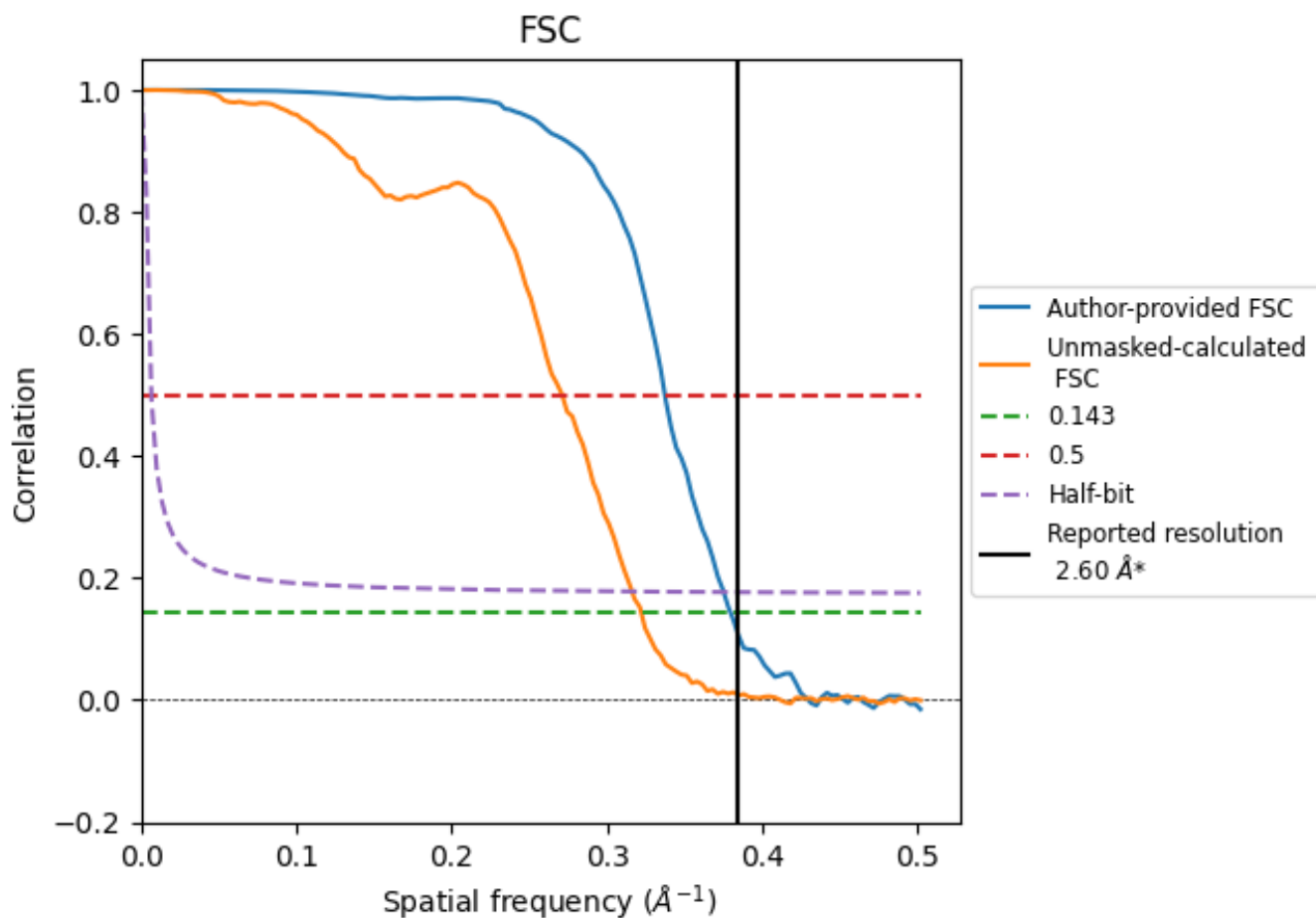


*Reported resolution corresponds to spatial frequency of 0.385 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.385\AA^{-1}

8.2 Resolution estimates [i](#)

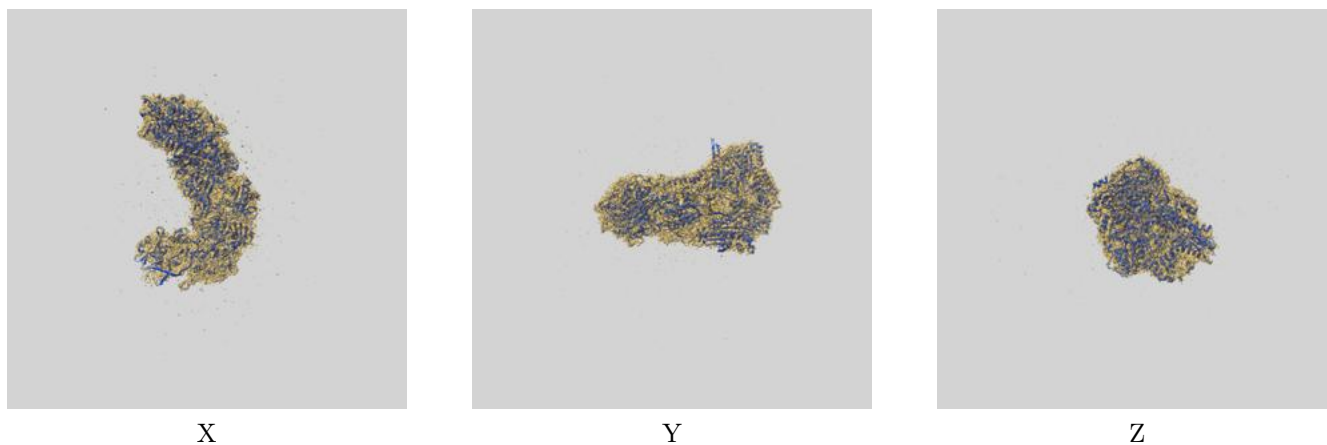
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.60	-	-
Author-provided FSC curve	2.63	2.96	2.66
Unmasked-calculated*	3.10	3.69	3.16

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.10 differs from the reported value 2.6 by more than 10 %

9 Map-model fit [i](#)

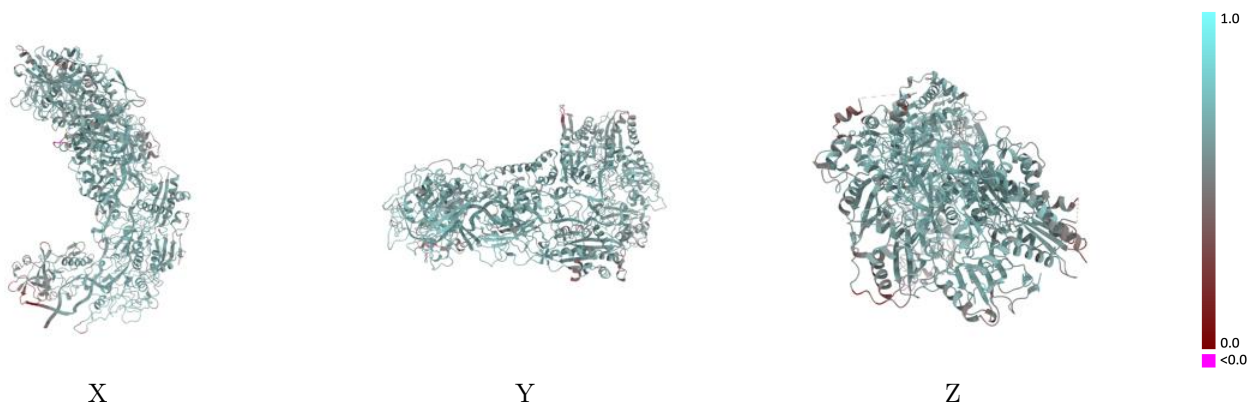
This section contains information regarding the fit between EMDB map EMD-66346 and PDB model 9WXH. Per-residue inclusion information can be found in section [3](#) on page [8](#).

9.1 Map-model overlay [i](#)



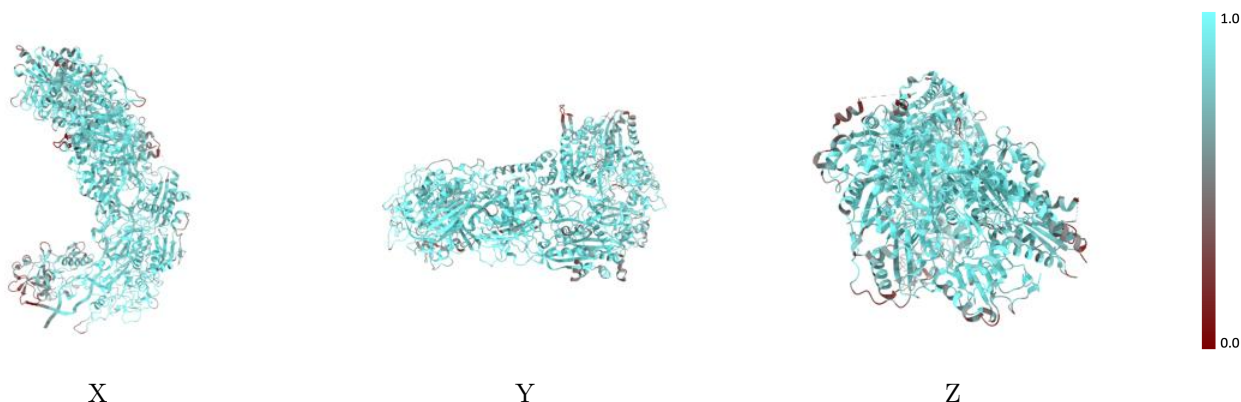
The images above show the 3D surface view of the map at the recommended contour level 0.111 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



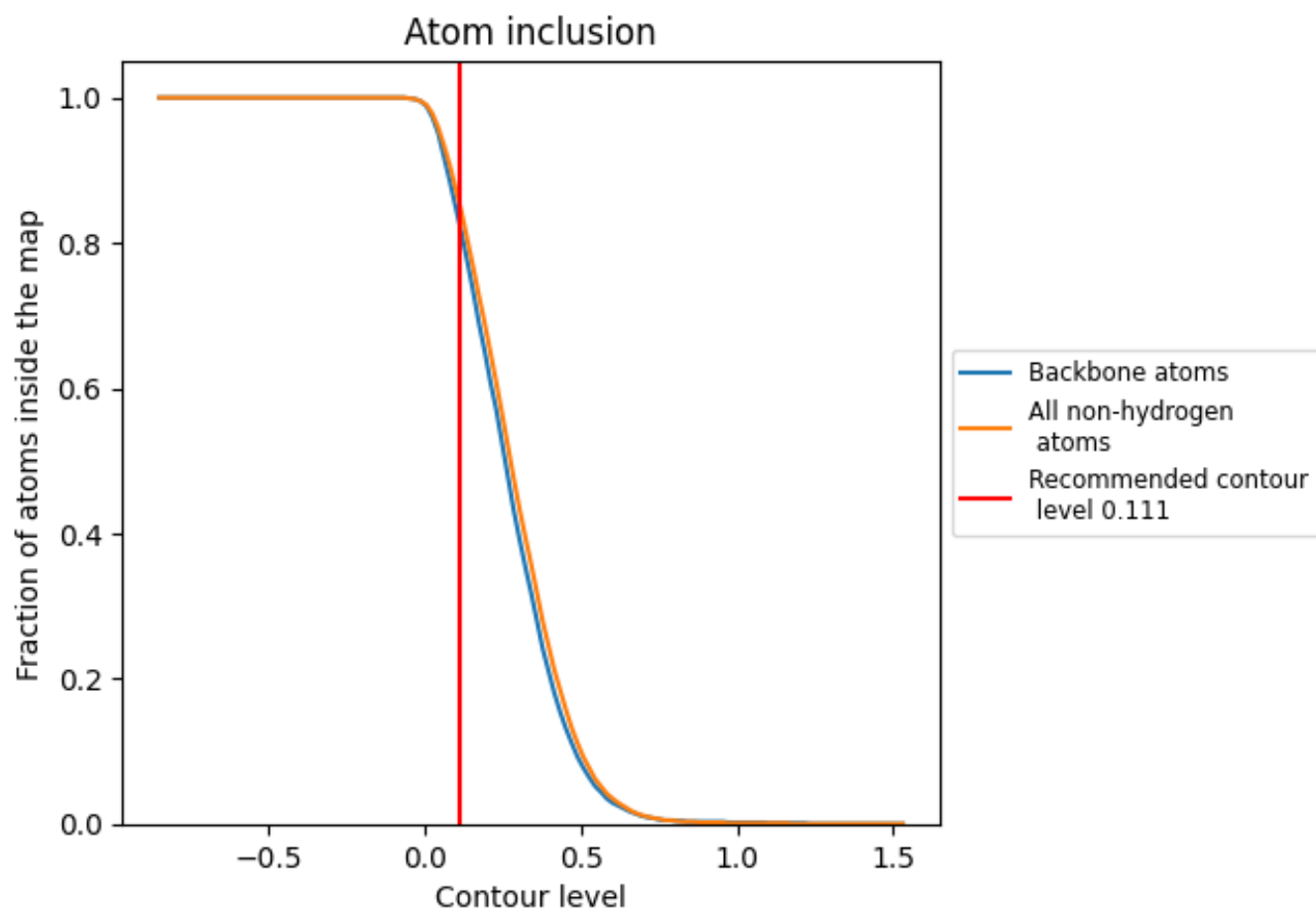
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.111).













9.4 Atom inclusion [i](#)



At the recommended contour level, 83% of all backbone atoms, 86% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary

The table lists the average atom inclusion at the recommended contour level (0.111) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8560	 0.6100
A	 0.9690	 0.6530
B	 0.8520	 0.6070
C	 0.8750	 0.6120
D	 0.8400	 0.6030
E	 0.9340	 0.6330

